

for 10/643,110

## EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	6455	precursor\$1 same nitride	US-PGPUB; USPAT; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/06/16 15:13
L2	292	precursor\$1 same nitride same halide\$1	US-PGPUB; USPAT; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/06/16 15:13
L3	125	2 and (amine or amide)	US-PGPUB; USPAT; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/06/16 15:13

06/16/2006 10643110 Everhart

=> d his full

(FILE 'HOME' ENTERED AT 10:21:17 ON 16 JUN 2006)

L1 FILE 'REGISTRY' ENTERED AT 10:21:30 ON 16 JUN 2006  
STRUCTURE UPLOADED  
D L1

L2 FILE 'MARPAT' ENTERED AT 10:23:09 ON 16 JUN 2006  
0 SEA ABB=ON PLU=ON US2005042888/PN

L3 FILE 'REGISTRY' ENTERED AT 11:04:30 ON 16 JUN 2006  
STRUCTURE UPLOADED  
D L3

L4 50 SEA SSS SAM L3  
L5 5629 SEA SSS FUL L3

FILE 'CAPLUS' ENTERED AT 11:06:28 ON 16 JUN 2006  
SET LINE 250  
SET DETAIL OFF  
E MOCVD+ALL/CT  
SET LINE LOGIN  
SET DETAIL LOGIN  
L6 285107 SEA ABB=ON PLU=ON MOCVD OR CVD OR VAPOR DEPOS? OR ATOMIC  
LAYER DEPOS? OR ALD OR ALCVD OR MONOLAYER OR ATOMIC LAYER  
EPITAR? OR BINARY REACTION OR SEQUENTIAL SURFACE REACT? OR  
(PULSED OR PULSING OR PULSE) (2W) (PRECUR?)  
L7 38 SEA ABB=ON PLU=ON L5 AND L6  
D IBIB ABS HITSTR HITIND 1-38

FILE 'STNGUIDE' ENTERED AT 11:13:56 ON 16 JUN 2006

L8 FILE 'REGISTRY' ENTERED AT 11:19:21 ON 16 JUN 2006  
STRUCTURE UPLOADED  
D L8  
L9 7 SEA SSS SAM L8  
L10 134 SEA SSS FUL L8

L11 FILE 'CAPLUS' ENTERED AT 11:20:28 ON 16 JUN 2006  
0 SEA ABB=ON PLU=ON L10 AND L6

FILE 'STNGUIDE' ENTERED AT 11:21:11 ON 16 JUN 2006

L12 FILE 'REGISTRY' ENTERED AT 11:22:27 ON 16 JUN 2006  
STRUCTURE UPLOADED  
D L12  
L13 15 SEA SSS SAM L12  
L14 341 SEA SSS FUL L12

L15 FILE 'CAPLUS' ENTERED AT 11:23:39 ON 16 JUN 2006  
1 SEA ABB=ON PLU=ON L14 AND L6  
L16 0 SEA ABB=ON PLU=ON L15 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:24:10 ON 16 JUN 2006

L17 FILE 'REGISTRY' ENTERED AT 11:25:35 ON 16 JUN 2006  
STRUCTURE UPLOADED

EIC 2800 MARY S. MIMS 272-5928

06/16/2006 10643110 Everhart

L18 9 SEA SSS SAM L17  
L19 811 SEA SSS FUL L17

FILE 'CAPLUS' ENTERED AT 11:27:06 ON 16 JUN 2006

L20 3 SEA ABB=ON PLU=ON L19 AND L6  
L21 0 SEA ABB=ON PLU=ON L20 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:28:37 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:30:10 ON 16 JUN 2006  
L22 STRUCTURE UPLOADED  
L23 3086 SEA SSS FUL L22

FILE 'CAPLUS' ENTERED AT 11:30:43 ON 16 JUN 2006

L24 28 SEA ABB=ON PLU=ON L23 AND L6  
L25 0 SEA ABB=ON PLU=ON L24 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:32:09 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:35:33 ON 16 JUN 2006  
L26 STRUCTURE UPLOADED  
D L26  
L27 0 SEA SSS SAM L26  
L28 0 SEA SSS FUL L26

FILE 'STNGUIDE' ENTERED AT 11:36:32 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:38:52 ON 16 JUN 2006  
L29 STRUCTURE UPLOADED  
D L29  
L30 7 SEA SSS SAM L29  
L31 274 SEA SSS FUL L29

FILE 'CAPLUS' ENTERED AT 11:40:02 ON 16 JUN 2006

L32 1 SEA ABB=ON PLU=ON L31 AND L6  
L33 0 SEA ABB=ON PLU=ON L32 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:40:36 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:42:34 ON 16 JUN 2006  
L34 STRUCTURE UPLOADED  
L35 0 SEA SSS SAM L34  
L36 0 SEA SSS FUL L34

FILE 'STNGUIDE' ENTERED AT 11:43:19 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:46:39 ON 16 JUN 2006  
L37 STRUCTURE UPLOADED  
L38 0 SEA SSS SAM L37  
L39 0 SEA SSS FUL L26

FILE 'STNGUIDE' ENTERED AT 11:48:39 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:52:57 ON 16 JUN 2006  
L40 STRUCTURE UPLOADED  
D L40  
L41 0 SEA SSS SAM L40  
L42 0 SEA SSS FUL L40  
L43 1 SEA ABB=ON PLU=ON 7440-33-7

EIC 2800 MARY S. MIMS 272-5928

06/16/2006 10643110 Everhart

D FIDE

FILE 'CAPLUS' ENTERED AT 11:57:16 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 13:40:35 ON 16 JUN 2006

L44 STRUCTURE UPLOADED

D L44

L45 50 SEA SSS SAM L44

L46 28399 SEA SSS FUL L44

FILE 'CAPLUS' ENTERED AT 13:42:39 ON 16 JUN 2006

L47 2336 SEA ABB=ON PLU=ON L46 AND L6

L48 349 SEA ABB=ON PLU=ON L47 AND ?PRECURSOR?

S SILICON NITRIDE/CN

FILE 'REGISTRY' ENTERED AT 13:45:14 ON 16 JUN 2006

L49 1 SEA ABB=ON PLU=ON SILICON NITRIDE/CN

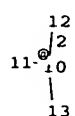
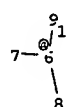
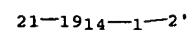
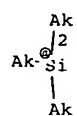
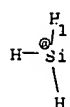
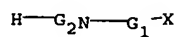
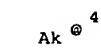
FILE 'CAPLUS' ENTERED AT 13:45:14 ON 16 JUN 2006

L50 78165 SEA ABB=ON PLU=ON L49

L51 23 SEA ABB=ON PLU=ON L48 AND L50

L52 22 SEA ABB=ON PLU=ON L51 NOT L7

D IBIB ABS HITSTR 1-22



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

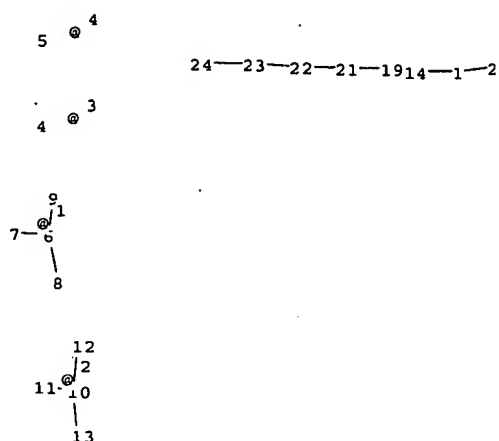
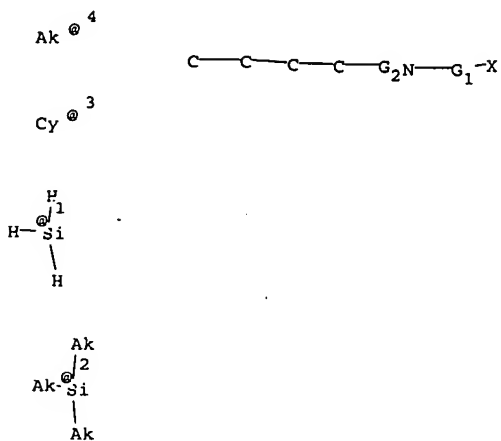
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS  
13:CLASS14:CLASS19:CLASS21:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21 22 23 24

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21 21-22 22-23 23-24

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

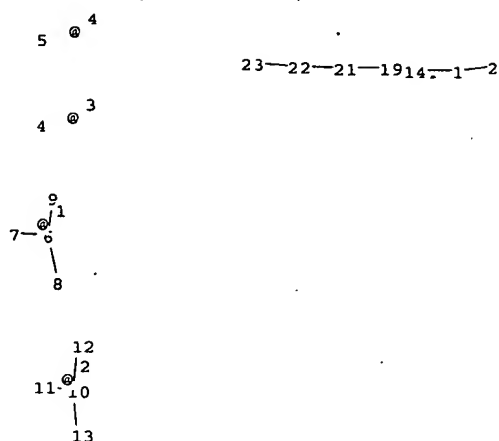
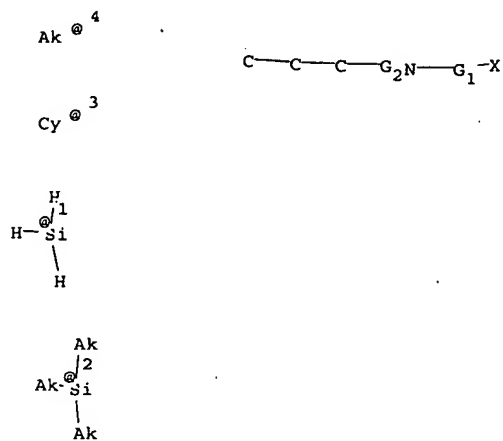
6-7 6-8 6-9 21-22 22-23 23-24

G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS23:CLASS24:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21 22 23

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21 21-22 22-23

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

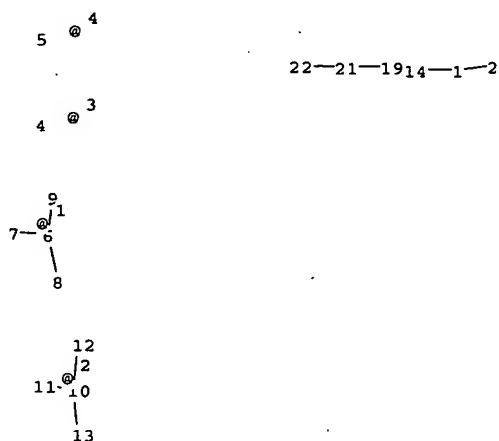
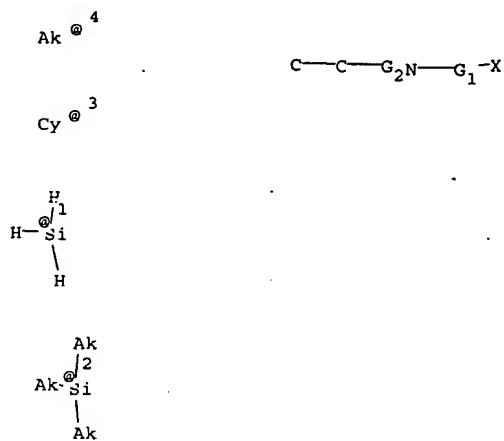
6-7 6-8 6-9 21-22 22-23

G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS23:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21 22

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21 21-22

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

6-7 6-8 6-9 21-22

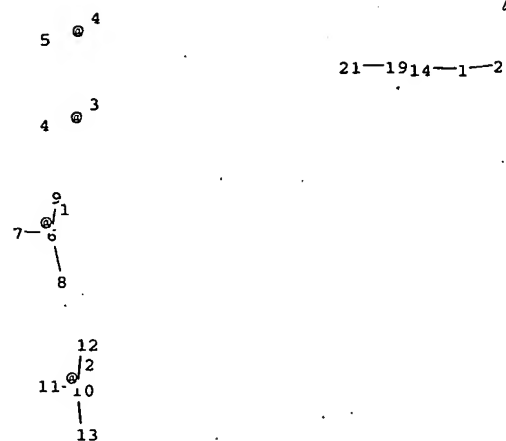
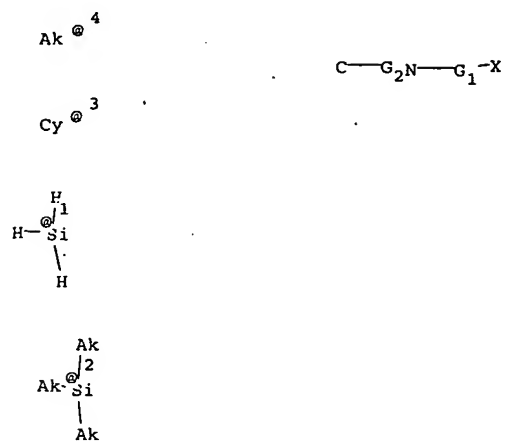
G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS  
13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS





chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS  
13:CLASS14:CLASS19:CLASS21:CLASS

1 2 4 5 6 7 8 9 10 11 12 13 14 19

21 22 23

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

21-22 21-23 22-23

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-23 22-23

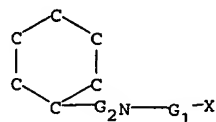
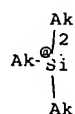
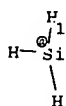
6-7 6-8 6-9

G2:[\*3],[\*4]

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS23:CLASS

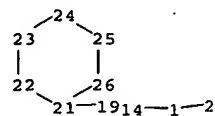
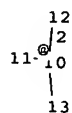
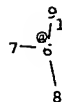
Ak @ 4

Cy @ 3



5 @ 4

4 @ 3



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23 24 25 26

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-26 22-23 23-24 24-25 25-26

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-26 22-23 23-24 24-25 25-26

exact bonds :

6-7 6-8 6-9

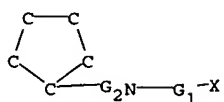
G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

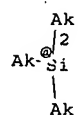
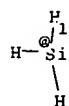
Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS  
13:CLASS14:CLASS19:CLASS21:Atom 22:CLASS23:CLASS24:CLASS25:Atom 26:Atom

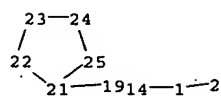
Ak @ 4



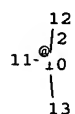
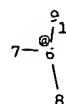
Cy @ 3



5 @ 4



4 @ 3



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23 24 25

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-25 22-23 23-24 24-25

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-25 22-23 23-24 24-25

exact bonds :

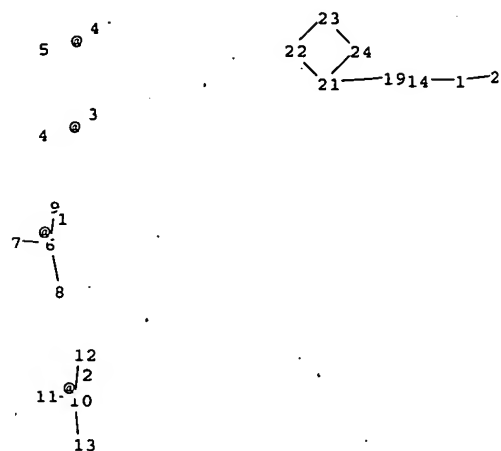
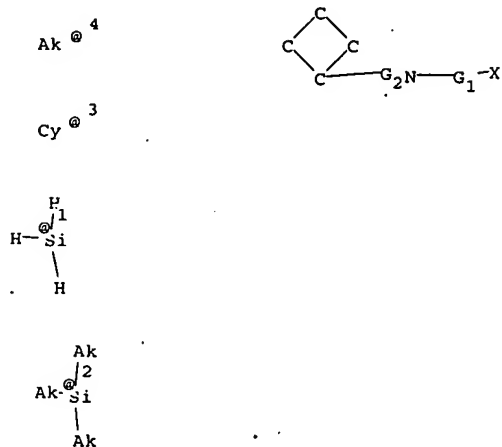
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS  
13:CLASS14:CLASS19:CLASS21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23 24

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-24 22-23 23-24

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-24 22-23 23-24

exact bonds :

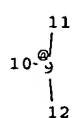
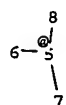
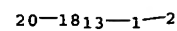
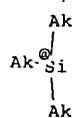
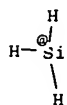
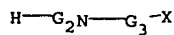
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[\*1],[\*2]

G2:[\*3],[\*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS  
13:CLASS14:CLASS19:CLASS21:Atom 22:Atom 23:Atom 24:CLASS



chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 18 20

chain bonds :

1-2 1-13 5-6 5-7 5-8 9-10 9-11 9-12 13-18 18-20

exact/norm bonds :

1-2 1-13 9-10 9-11 9-12 13-18 18-20

exact bonds :

5-6 5-7 5-8

G2:[\*1],[\*2]

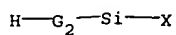
G3:Ta,Ti

Match level :

1:CLASS2:CLASS3:Atom 4:CLASS5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS  
12:CLASS13:CLASS18:CLASS20:CLASS

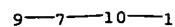
Ak @<sup>2</sup>

Cy @<sup>1</sup>



4 @<sup>2</sup>

3 @<sup>1</sup>



chain nodes :

1 3 4 7 9 10

chain bonds :

1-10 7-9 7-10

exact/norm bonds :

7-9 7-10

exact bonds :

1-10

G1:Si,B,Al,Nb,Ta,Ti,W

G2:[\*1],[\*2]

Match level :

1:CLASS3:Atom 4:CLASS7:CLASS9:CLASS10:CLASS